



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 3, 2018 – 10:57 PM EST

PDB ID : 5WS8
Title : Pyruvate kinase (PYK) from Mycobacterium tuberculosis in complex with Oxalate
Authors : Zhong, W.; Cai, Q.; El Sahili, A.; Lescar, J.; Dedon, P.C.
Deposited on : 2016-12-06
Resolution : 2.62 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030736
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030736

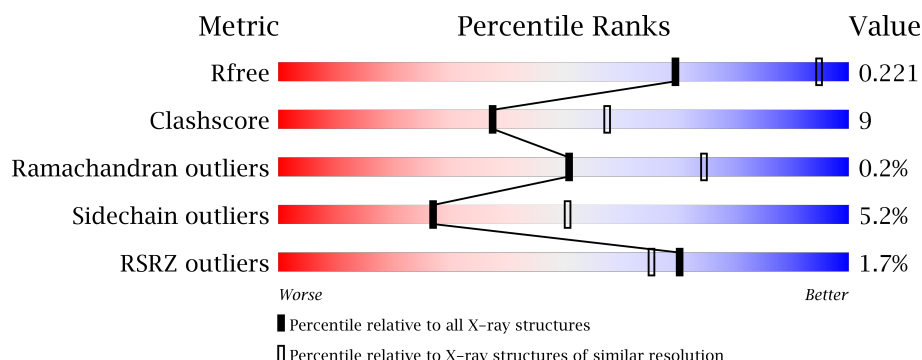
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.62 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	2983 (2.64-2.60)
Clashscore	112137	3351 (2.64-2.60)
Ramachandran outliers	110173	3298 (2.64-2.60)
Sidechain outliers	110143	3298 (2.64-2.60)
RSRZ outliers	101464	2992 (2.64-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	475	<div> <div>0.2%</div> <div>78%</div> <div>18%</div> <div>..</div> </div>
1	B	475	<div> <div>2%</div> <div>78%</div> <div>18%</div> <div>..</div> </div>
1	C	475	<div> <div>2%</div> <div>78%</div> <div>18%</div> <div>..</div> </div>
1	D	475	<div> <div>0.2%</div> <div>80%</div> <div>16%</div> <div>..</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14566 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Pyruvate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	1	0
			3511	2185	635	673	18			
1	B	464	Total	C	N	O	S	0	1	0
			3511	2185	635	673	18			
1	C	464	Total	C	N	O	S	0	1	0
			3511	2185	635	673	18			
1	D	464	Total	C	N	O	S	0	1	0
			3511	2185	635	673	18			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P9WKE5
A	-1	GLY	-	expression tag	UNP P9WKE5
A	0	HIS	-	expression tag	UNP P9WKE5
B	-2	GLY	-	expression tag	UNP P9WKE5
B	-1	GLY	-	expression tag	UNP P9WKE5
B	0	HIS	-	expression tag	UNP P9WKE5
C	-2	GLY	-	expression tag	UNP P9WKE5
C	-1	GLY	-	expression tag	UNP P9WKE5
C	0	HIS	-	expression tag	UNP P9WKE5
D	-2	GLY	-	expression tag	UNP P9WKE5
D	-1	GLY	-	expression tag	UNP P9WKE5
D	0	HIS	-	expression tag	UNP P9WKE5

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

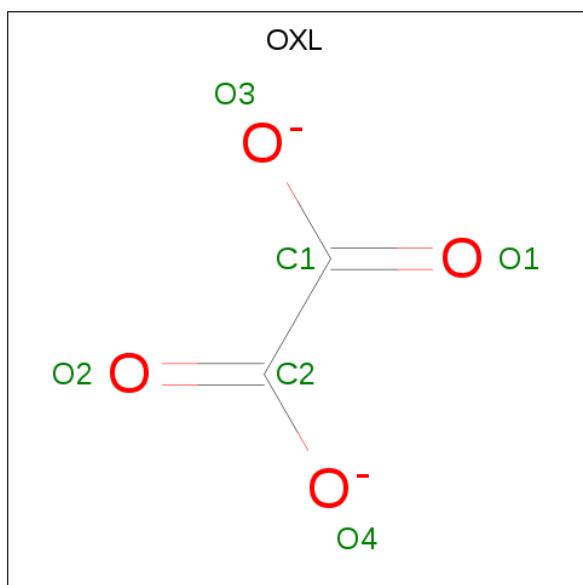
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Mg	0	0
			1	1		
2	C	1	Total	Mg	0	0
			1	1		

- Molecule 3 is OXALATE ION (three-letter code: OXL) (formula: C_2O_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	2	4		
3	B	1	Total	C	O	0	0
			6	2	4		
3	C	1	Total	C	O	0	0
			6	2	4		
3	D	1	Total	C	O	0	0
			6	2	4		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	130	Total	O	0	0
			130	130		
4	B	110	Total	O	0	0
			110	110		
4	C	123	Total	O	0	0
			123	123		

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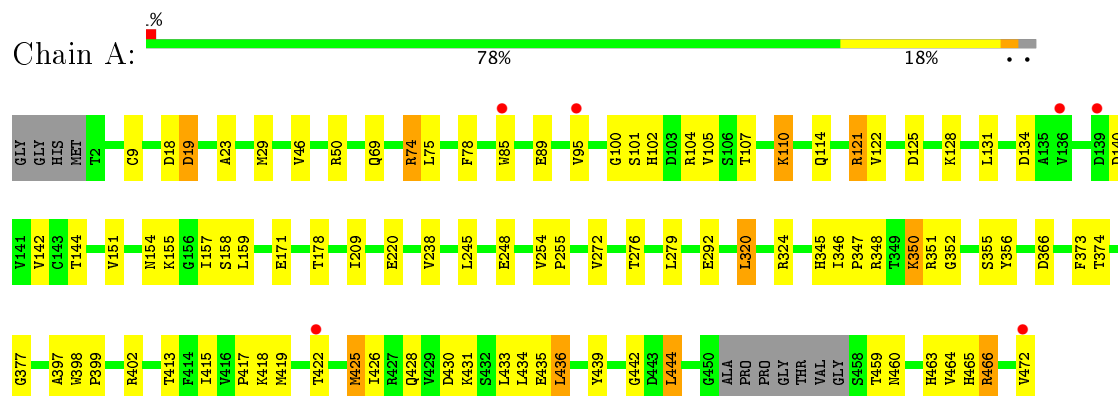
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	131	Total	O	0	0
			131	131		

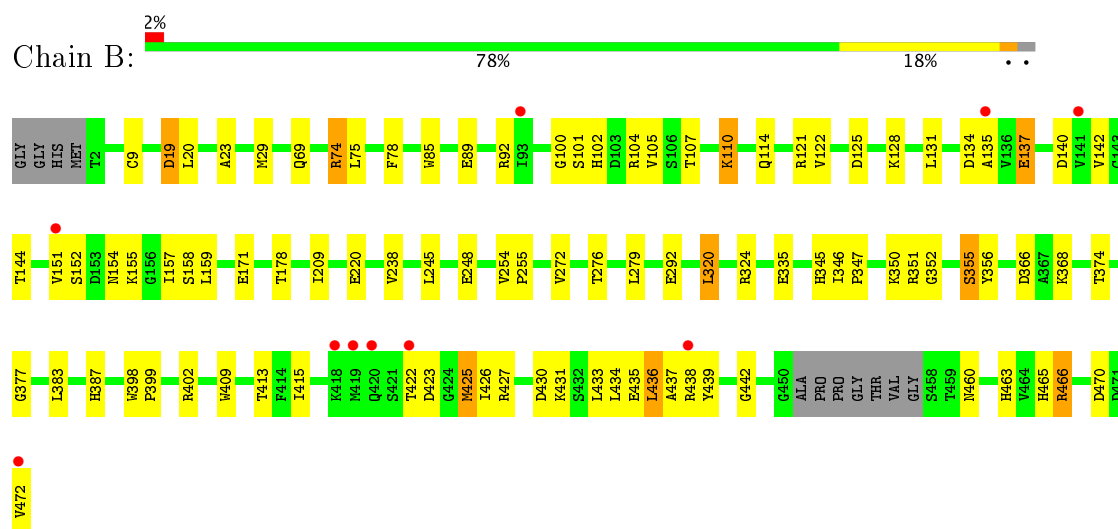
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

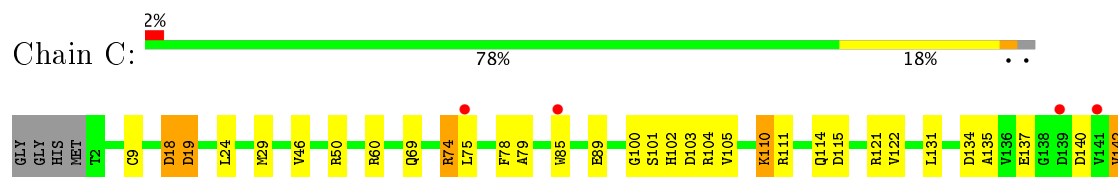
• Molecule 1: Pyruvate kinase

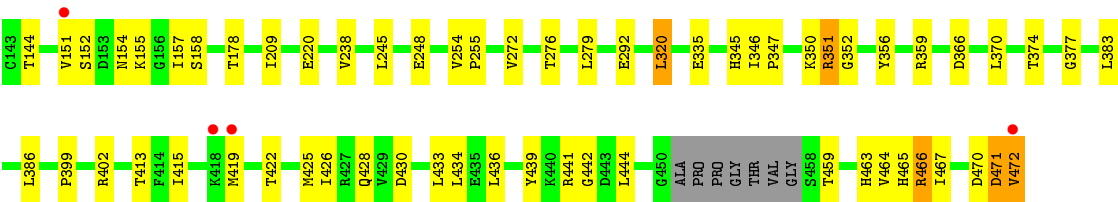


• Molecule 1: Pyruvate kinase

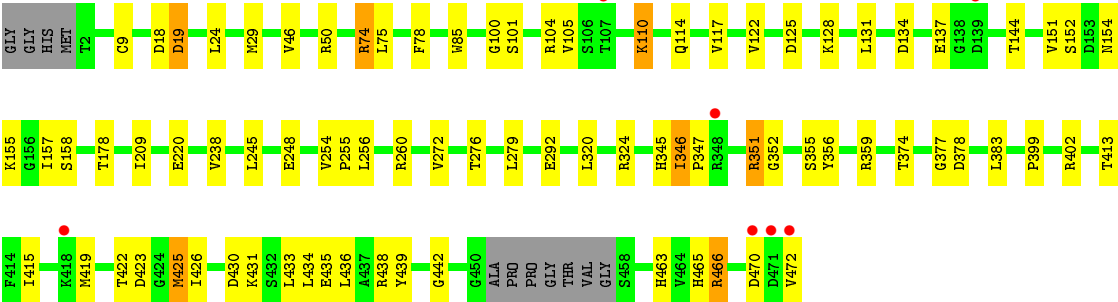
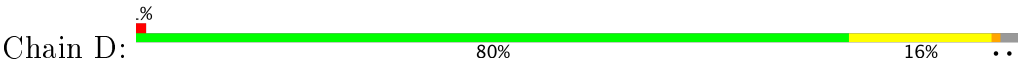


• Molecule 1: Pyruvate kinase





● Molecule 1: Pyruvate kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	127.01Å 127.01Å 144.65Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	63.51 – 2.62 44.16 – 2.62	Depositor EDS
% Data completeness (in resolution range)	100.0 (63.51-2.62) 100.0 (44.16-2.62)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.173 , 0.218 0.181 , 0.221	Depositor DCC
R_{free} test set	4011 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	49.6	Xtriage
Anisotropy	0.703	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.477 for -h,-k,l 0.477 for h,-h-k,-l 0.480 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14566	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: OXL, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.60	0/3558	0.64	0/4829
1	B	0.60	0/3558	0.63	0/4829
1	C	0.61	0/3558	0.65	0/4829
1	D	0.60	0/3558	0.64	0/4829
All	All	0.60	0/14232	0.64	0/19316

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3511	0	3573	73	0
1	B	3511	0	3573	57	0
1	C	3511	0	3573	64	0
1	D	3511	0	3573	56	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	6	0	0	0	0
3	C	6	0	0	0	0
3	D	6	0	0	0	0
4	A	130	0	0	5	0
4	B	110	0	0	3	0
4	C	123	0	0	8	0
4	D	131	0	0	6	0
All	All	14566	0	14292	249	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (249) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:351:ARG:NH2	1:D:378:ASP:OD1	1.82	1.12
1:C:430:ASP:O	1:C:434:LEU:HD12	1.58	1.03
1:B:320:LEU:HD23	1:B:324:ARG:NH1	1.82	0.94
1:D:347:PRO:HG3	1:D:356:TYR:CE2	2.04	0.91
1:D:125:ASP:OD2	1:D:128:LYS:HD2	1.70	0.91
1:A:122:VAL:HG22	1:A:159:LEU:HD11	1.52	0.89
1:B:423:ASP:O	1:B:427:ARG:HG3	1.76	0.85
1:B:92:ARG:NH1	1:B:137:GLU:OE2	2.10	0.84
1:A:95:VAL:HG13	1:A:110:LYS:HA	1.58	0.83
1:A:320:LEU:HD23	1:A:324:ARG:NH1	1.94	0.81
1:C:135:ALA:O	1:C:142:VAL:HG23	1.82	0.79
1:A:69:GLN:OE1	1:A:74:ARG:NH1	2.17	0.77
1:A:418:LYS:H	1:A:418:LYS:HD2	1.50	0.77
1:C:85:TRP:CE3	1:C:151:VAL:HG11	2.21	0.76
1:C:347:PRO:HG3	1:C:356:TYR:CE2	2.21	0.75
1:B:85:TRP:CE3	1:B:151:VAL:HG11	2.22	0.75
1:A:347:PRO:HG3	1:A:356:TYR:CE2	2.21	0.75
1:D:85:TRP:CE3	1:D:151:VAL:HG11	2.22	0.75
1:A:85:TRP:CE3	1:A:151:VAL:HG11	2.22	0.75
1:B:347:PRO:HG3	1:B:356:TYR:CE2	2.22	0.74
1:B:436:LEU:N	1:B:436:LEU:HD12	2.03	0.73
1:A:122:VAL:HG22	1:A:159:LEU:CD1	2.17	0.73
1:C:18:ASP:HA	1:C:50:ARG:HH12	1.54	0.73
1:A:436:LEU:HD12	1:A:436:LEU:N	2.05	0.71
1:A:402:ARG:HG3	1:A:413:THR:HB	1.72	0.71
1:C:402:ARG:HG3	1:C:413:THR:HB	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:402:ARG:HG3	1:D:413:THR:HB	1.73	0.70
1:B:402:ARG:HG3	1:B:413:THR:HB	1.75	0.69
1:A:419:MET:SD	1:A:428:GLN:HG2	2.33	0.68
1:D:74:ARG:NH1	1:D:154:ASN:O	2.26	0.68
1:B:426:ILE:O	1:B:430:ASP:OD2	2.13	0.66
1:D:426:ILE:O	1:D:430:ASP:OD2	2.13	0.66
1:A:158:SER:OG	1:A:248:GLU:OE1	2.14	0.66
1:D:158:SER:OG	1:D:248:GLU:OE1	2.14	0.66
1:B:158:SER:OG	1:B:248:GLU:OE1	2.14	0.66
1:C:426:ILE:O	1:C:430:ASP:OD2	2.13	0.66
1:C:436:LEU:N	1:C:436:LEU:HD12	2.10	0.66
1:C:158:SER:OG	1:C:248:GLU:OE1	2.14	0.66
1:A:426:ILE:O	1:A:430:ASP:OD2	2.14	0.66
1:B:134:ASP:OD2	1:B:144:THR:OG1	2.15	0.65
1:D:125:ASP:OD2	1:D:128:LYS:CD	2.44	0.65
1:A:178:THR:OG1	1:A:209:ILE:HD11	1.97	0.65
1:A:134:ASP:OD2	1:A:144:THR:OG1	2.16	0.64
1:C:430:ASP:O	1:C:434:LEU:CD1	2.40	0.64
1:C:134:ASP:OD2	1:C:144:THR:OG1	2.15	0.64
1:D:134:ASP:OD2	1:D:144:THR:OG1	2.15	0.64
1:D:436:LEU:N	1:D:436:LEU:HD12	2.12	0.64
1:C:140:ASP:HB2	4:C:621:HOH:O	1.97	0.63
1:A:238:VAL:HG23	1:A:272:VAL:HG23	1.81	0.62
1:C:178:THR:OG1	1:C:209:ILE:HD11	1.99	0.62
1:C:104:ARG:NH2	4:C:601:HOH:O	2.32	0.62
1:B:238:VAL:HG23	1:B:272:VAL:HG23	1.82	0.61
1:C:238:VAL:HG23	1:C:272:VAL:HG23	1.81	0.61
1:D:260:ARG:HD3	4:D:675:HOH:O	2.01	0.61
1:D:178:THR:OG1	1:D:209:ILE:HD11	2.01	0.61
1:D:238:VAL:HG23	1:D:272:VAL:HG23	1.81	0.61
1:B:69:GLN:OE1	1:B:74:ARG:NH1	2.34	0.61
1:B:178:THR:OG1	1:B:209:ILE:HD11	2.01	0.61
1:C:441:ARG:HA	1:C:467:ILE:HG22	1.81	0.61
1:D:346:ILE:H	1:D:346:ILE:HD12	1.66	0.60
1:C:430:ASP:C	1:C:434:LEU:HD12	2.22	0.60
1:A:140:ASP:HB2	4:A:628:HOH:O	2.03	0.59
1:C:78:PHE:CZ	1:C:105:VAL:HG21	2.38	0.59
1:B:78:PHE:CZ	1:B:105:VAL:HG21	2.38	0.58
1:C:69:GLN:OE1	1:C:74:ARG:NH2	2.36	0.58
1:A:78:PHE:CZ	1:A:105:VAL:HG21	2.39	0.58
1:D:347:PRO:CG	1:D:356:TYR:CE2	2.83	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:29:MET:O	1:C:60:ARG:NH2	2.37	0.57
1:B:125:ASP:OD2	1:B:128:LYS:HG3	2.04	0.57
1:D:78:PHE:CZ	1:D:105:VAL:HG21	2.39	0.57
1:A:419:MET:CE	1:A:428:GLN:HG2	2.36	0.56
1:D:78:PHE:HA	1:D:100:GLY:O	2.07	0.55
1:D:104:ARG:NH2	4:D:602:HOH:O	2.39	0.55
1:A:422:THR:HA	1:A:425:MET:HG3	1.88	0.55
1:A:399:PRO:HA	1:A:415:ILE:HD12	1.87	0.55
1:A:95:VAL:HG13	1:A:110:LYS:HD2	1.89	0.55
1:D:347:PRO:CG	1:D:356:TYR:CD2	2.90	0.55
1:A:419:MET:HE1	1:A:428:GLN:HG2	1.89	0.55
1:C:399:PRO:HA	1:C:415:ILE:HD12	1.88	0.55
1:D:355:SER:HB3	1:D:383:LEU:HD13	1.87	0.55
1:A:151:VAL:HG23	1:A:151:VAL:O	2.07	0.54
1:C:370:LEU:HD11	1:C:383:LEU:HD22	1.89	0.54
1:A:122:VAL:HB	1:A:131:LEU:HB2	1.89	0.54
1:B:347:PRO:CG	1:B:356:TYR:CE2	2.89	0.54
1:C:121:ARG:HD2	4:D:661:HOH:O	2.08	0.54
1:A:460:ASN:ND2	1:C:471:ASP:HB2	2.23	0.54
1:A:347:PRO:CG	1:A:356:TYR:CE2	2.90	0.54
1:B:347:PRO:HG2	1:B:356:TYR:CD2	2.42	0.54
1:C:347:PRO:HG2	1:C:356:TYR:CD2	2.43	0.54
1:B:351:ARG:O	1:B:355:SER:OG	2.25	0.54
1:B:399:PRO:HA	1:B:415:ILE:HD12	1.89	0.53
1:C:422:THR:HA	1:C:425:MET:HG3	1.89	0.53
1:C:19:ASP:HA	4:C:622:HOH:O	2.07	0.53
1:D:347:PRO:HG3	1:D:356:TYR:CD2	2.43	0.53
1:D:399:PRO:HA	1:D:415:ILE:HD12	1.90	0.53
1:A:78:PHE:HA	1:A:100:GLY:O	2.09	0.53
1:A:19:ASP:HB2	4:A:610:HOH:O	2.09	0.53
1:C:347:PRO:CG	1:C:356:TYR:CE2	2.90	0.52
1:A:347:PRO:HG2	1:A:356:TYR:CD2	2.44	0.52
1:D:75:LEU:O	1:D:154:ASN:HA	2.10	0.52
1:B:436:LEU:N	1:B:436:LEU:CD1	2.72	0.52
1:A:347:PRO:CG	1:A:356:TYR:CD2	2.93	0.52
1:B:78:PHE:HA	1:B:100:GLY:O	2.10	0.52
1:C:75:LEU:O	1:C:154:ASN:HA	2.10	0.52
1:D:422:THR:HA	1:D:425:MET:HG3	1.91	0.52
1:B:347:PRO:CG	1:B:356:TYR:CD2	2.94	0.51
1:A:436:LEU:CD1	1:A:436:LEU:N	2.73	0.51
1:A:75:LEU:O	1:A:154:ASN:HA	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:422:THR:HA	1:B:425:MET:HG3	1.92	0.51
1:D:46:VAL:O	1:D:50:ARG:HG3	2.11	0.51
1:A:125:ASP:OD2	1:A:128:LYS:HG3	2.10	0.51
1:A:95:VAL:HG11	1:A:110:LYS:O	2.10	0.51
1:B:75:LEU:O	1:B:154:ASN:HA	2.10	0.51
1:D:122:VAL:HB	1:D:131:LEU:HB2	1.91	0.51
1:D:18:ASP:HA	1:D:50:ARG:HH22	1.74	0.51
1:C:347:PRO:CG	1:C:356:TYR:CD2	2.93	0.51
1:C:79:ALA:HB2	4:C:604:HOH:O	2.10	0.51
1:B:151:VAL:HG23	1:B:151:VAL:O	2.10	0.51
1:C:151:VAL:O	1:C:151:VAL:HG23	2.10	0.50
1:C:436:LEU:CD1	1:C:436:LEU:N	2.75	0.50
1:D:151:VAL:O	1:D:151:VAL:HG23	2.11	0.50
1:C:89:GLU:OE2	1:C:102:HIS:CE1	2.64	0.50
1:A:23:ALA:HB1	1:A:320:LEU:HD11	1.94	0.50
1:C:78:PHE:HA	1:C:100:GLY:O	2.12	0.49
1:D:78:PHE:HB3	4:D:687:HOH:O	2.11	0.49
1:A:373:PHE:HE2	1:A:419:MET:CE	2.26	0.49
1:B:352:GLY:HA3	4:B:644:HOH:O	2.13	0.49
1:C:352:GLY:HA3	4:C:666:HOH:O	2.13	0.49
1:D:436:LEU:CD1	1:D:436:LEU:N	2.76	0.49
1:D:430:ASP:O	1:D:434:LEU:HD12	2.14	0.48
1:A:426:ILE:HD13	1:A:472:VAL:HG22	1.95	0.48
1:B:135:ALA:O	1:B:142:VAL:HG23	2.14	0.48
1:C:374:THR:HG21	1:C:377:GLY:HA2	1.95	0.48
1:C:101:SER:O	1:C:104:ARG:O	2.32	0.48
1:B:101:SER:O	1:B:104:ARG:O	2.32	0.48
1:D:374:THR:HG21	1:D:377:GLY:HA2	1.96	0.48
1:D:101:SER:O	1:D:104:ARG:O	2.31	0.48
1:B:19:ASP:HA	4:B:619:HOH:O	2.14	0.48
1:D:433:LEU:O	1:D:439:TYR:HB2	2.14	0.48
1:D:347:PRO:HG2	1:D:356:TYR:CD2	2.48	0.47
1:A:397:ALA:HB1	1:A:418:LYS:HE3	1.95	0.47
1:B:426:ILE:HD13	1:B:472:VAL:HG22	1.96	0.47
1:C:350:LYS:NZ	1:C:459:THR:O	2.47	0.47
1:D:426:ILE:HD13	1:D:472:VAL:HG22	1.96	0.47
1:B:122:VAL:HB	1:B:131:LEU:HB2	1.96	0.47
1:C:89:GLU:OE2	1:C:102:HIS:NE2	2.47	0.47
1:C:419:MET:CE	1:C:428:GLN:HG2	2.44	0.47
1:B:430:ASP:O	1:B:434:LEU:HD12	2.14	0.47
1:A:374:THR:HG21	1:A:377:GLY:HA2	1.95	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:LYS:HA	1:A:110:LYS:HD2	1.66	0.46
1:A:121:ARG:HG2	4:B:662:HOH:O	2.14	0.46
1:D:320:LEU:HD23	1:D:324:ARG:HH12	1.80	0.46
1:A:430:ASP:O	1:A:434:LEU:HD12	2.14	0.46
1:B:374:THR:HG21	1:B:377:GLY:HA2	1.96	0.46
1:A:19:ASP:HA	4:A:616:HOH:O	2.15	0.46
1:D:9:CYS:HB2	1:D:29:MET:SD	2.56	0.46
1:B:89:GLU:OE1	1:B:102:HIS:NE2	2.49	0.46
1:B:171:GLU:OE2	1:B:171:GLU:HA	2.16	0.46
1:D:351:ARG:CZ	1:D:378:ASP:OD1	2.59	0.46
1:C:9:CYS:HB2	1:C:29:MET:SD	2.56	0.46
1:C:370:LEU:HD11	1:C:383:LEU:CD2	2.45	0.46
1:A:463:HIS:CD2	1:A:465:HIS:HB2	2.50	0.46
1:B:426:ILE:HD13	1:B:472:VAL:HG13	1.98	0.46
1:A:442:GLY:O	1:A:466:ARG:HG3	2.16	0.45
1:C:433:LEU:O	1:C:439:TYR:HB2	2.16	0.45
1:D:110:LYS:HD2	1:D:110:LYS:HA	1.70	0.45
1:B:431:LYS:O	1:B:435:GLU:HG3	2.16	0.45
4:C:708:HOH:O	1:D:256:LEU:HD12	2.17	0.45
1:A:433:LEU:O	1:A:439:TYR:HB2	2.17	0.45
1:C:279:LEU:HD23	1:C:292:GLU:HB3	1.99	0.45
1:D:426:ILE:HD13	1:D:472:VAL:HG13	1.99	0.45
1:B:463:HIS:CD2	1:B:465:HIS:HB2	2.51	0.45
1:D:431:LYS:O	1:D:435:GLU:HG3	2.16	0.45
1:C:463:HIS:CD2	1:C:465:HIS:HB2	2.52	0.44
1:D:279:LEU:HD23	1:D:292:GLU:HB3	1.98	0.44
1:A:89:GLU:OE1	1:A:102:HIS:CE1	2.70	0.44
1:A:347:PRO:HG3	1:A:356:TYR:CD2	2.52	0.44
1:C:110:LYS:HA	1:C:110:LYS:HD2	1.68	0.44
1:C:122:VAL:HB	1:C:131:LEU:HB2	1.98	0.44
1:C:351:ARG:HD3	4:C:689:HOH:O	2.17	0.44
1:A:431:LYS:O	1:A:435:GLU:HG3	2.18	0.44
1:B:101:SER:N	1:B:104:ARG:O	2.51	0.44
1:C:442:GLY:O	1:C:466:ARG:HG3	2.18	0.44
1:A:9:CYS:HB2	1:A:29:MET:SD	2.57	0.44
1:B:9:CYS:HB2	1:B:29:MET:SD	2.58	0.44
1:B:433:LEU:O	1:B:439:TYR:HB2	2.18	0.44
1:B:254:VAL:N	1:B:255:PRO:CD	2.81	0.44
1:B:279:LEU:HD23	1:B:292:GLU:HB3	1.99	0.44
1:C:101:SER:N	1:C:104:ARG:O	2.51	0.44
1:C:24:LEU:HD23	1:C:320:LEU:HD12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:442:GLY:O	1:B:466:ARG:HG3	2.18	0.43
1:A:426:ILE:HD13	1:A:472:VAL:HG13	2.00	0.43
1:D:320:LEU:HD12	1:D:320:LEU:HA	1.63	0.43
1:A:254:VAL:N	1:A:255:PRO:CD	2.82	0.43
1:A:350:LYS:HE2	1:A:459:THR:O	2.18	0.43
1:C:346:ILE:HD12	1:C:347:PRO:HD3	2.01	0.43
1:A:46:VAL:O	1:A:50:ARG:HG3	2.18	0.43
1:A:78:PHE:HB3	4:A:687:HOH:O	2.19	0.43
1:B:423:ASP:OD1	1:B:472:VAL:HG11	2.18	0.43
1:B:110:LYS:HD2	1:B:110:LYS:HA	1.69	0.43
1:B:346:ILE:HD12	1:B:346:ILE:HA	1.91	0.43
1:B:346:ILE:HA	1:B:347:PRO:HD3	1.87	0.42
1:D:125:ASP:O	1:D:128:LYS:HG3	2.19	0.42
1:A:95:VAL:CG1	1:A:110:LYS:HA	2.40	0.42
1:A:279:LEU:HD23	1:A:292:GLU:HB3	2.01	0.42
1:C:347:PRO:HG3	1:C:356:TYR:CD2	2.53	0.42
1:D:463:HIS:CD2	1:D:465:HIS:HB2	2.54	0.42
1:A:171:GLU:HA	1:A:171:GLU:OE2	2.17	0.42
1:A:346:ILE:HA	1:A:347:PRO:HD3	1.87	0.42
1:B:437:ALA:O	1:B:438:ARG:HB2	2.19	0.42
1:C:444:LEU:HD11	1:C:464:VAL:CG1	2.49	0.42
1:D:423:ASP:OD1	1:D:472:VAL:HG11	2.19	0.42
1:A:220:GLU:HA	1:A:245:LEU:HD13	2.00	0.42
1:A:348:ARG:CZ	1:A:348:ARG:HB2	2.49	0.42
1:B:220:GLU:HA	1:B:245:LEU:HD13	2.00	0.42
1:B:398:TRP:HA	1:B:399:PRO:HD3	1.85	0.42
1:B:350:LYS:HE3	1:B:460:ASN:HB3	2.02	0.42
1:C:335:GLU:HB2	4:C:633:HOH:O	2.19	0.42
1:D:254:VAL:N	1:D:255:PRO:CD	2.82	0.42
1:A:351:ARG:O	1:A:355:SER:OG	2.34	0.42
1:D:442:GLY:O	1:D:466:ARG:HG3	2.20	0.42
1:A:134:ASP:HB2	1:A:142:VAL:HG13	2.02	0.42
1:B:107:THR:O	1:B:110:LYS:NZ	2.47	0.42
1:C:346:ILE:HA	1:C:347:PRO:HD3	1.86	0.42
1:D:352:GLY:HA3	4:D:673:HOH:O	2.20	0.42
1:A:107:THR:O	1:A:110:LYS:NZ	2.46	0.41
1:A:352:GLY:HA3	4:A:690:HOH:O	2.20	0.41
1:C:254:VAL:N	1:C:255:PRO:CD	2.82	0.41
1:A:463:HIS:HD2	1:A:465:HIS:HB2	1.85	0.41
1:C:419:MET:HE3	1:C:428:GLN:HG2	2.02	0.41
1:A:23:ALA:CB	1:A:320:LEU:HD11	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:398:TRP:HA	1:A:399:PRO:HD3	1.84	0.41
1:D:101:SER:N	1:D:104:ARG:O	2.53	0.41
1:D:220:GLU:HA	1:D:245:LEU:HD13	2.00	0.41
1:A:444:LEU:HD21	1:A:464:VAL:CG1	2.51	0.41
1:D:19:ASP:HA	4:D:613:HOH:O	2.19	0.41
1:B:23:ALA:CB	1:B:320:LEU:HD11	2.51	0.41
1:C:220:GLU:HA	1:C:245:LEU:HD13	2.01	0.41
1:A:417:PRO:O	1:A:419:MET:SD	2.78	0.41
1:C:426:ILE:HD13	1:C:472:VAL:HB	2.03	0.41
1:C:111:ARG:O	1:C:115:ASP:HB2	2.21	0.40
1:C:19:ASP:H	1:C:50:ARG:NH1	2.19	0.40
1:A:121:ARG:O	1:A:159:LEU:HD13	2.21	0.40
1:B:320:LEU:HD23	1:B:324:ARG:CZ	2.48	0.40
1:C:46:VAL:O	1:C:50:ARG:HG3	2.21	0.40
1:D:24:LEU:HD23	1:D:320:LEU:HD12	2.02	0.40
1:A:101:SER:N	1:A:104:ARG:O	2.55	0.40
1:B:346:ILE:HD12	1:B:347:PRO:HD3	2.04	0.40
1:B:387:HIS:HA	1:B:409:TRP:CZ3	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	461/475 (97%)	445 (96%)	15 (3%)	1 (0%)	51	74
1	B	461/475 (97%)	446 (97%)	14 (3%)	1 (0%)	51	74
1	C	461/475 (97%)	444 (96%)	16 (4%)	1 (0%)	51	74
1	D	461/475 (97%)	447 (97%)	13 (3%)	1 (0%)	51	74
All	All	1844/1900 (97%)	1782 (97%)	58 (3%)	4 (0%)	51	74

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	276	THR
1	A	276	THR
1	B	276	THR
1	D	276	THR

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	376/381 (99%)	360 (96%)	16 (4%)	33	60
1	B	376/381 (99%)	353 (94%)	23 (6%)	22	42
1	C	376/381 (99%)	355 (94%)	21 (6%)	25	46
1	D	376/381 (99%)	358 (95%)	18 (5%)	30	54
All	All	1504/1524 (99%)	1426 (95%)	78 (5%)	27	51

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ASP
1	A	19	ASP
1	A	74	ARG
1	A	110	LYS
1	A	114	GLN
1	A	121	ARG
1	A	155	LYS
1	A	157	ILE
1	A	320	LEU
1	A	345	HIS
1	A	350	LYS
1	A	366	ASP
1	A	425	MET
1	A	436	LEU
1	A	444	LEU
1	A	466	ARG
1	B	19	ASP
1	B	20	LEU

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Mol	Chain	Res	Type
1	B	74	ARG
1	B	110	LYS
1	B	114	GLN
1	B	121	ARG
1	B	137	GLU
1	B	140	ASP
1	B	152	SER
1	B	155	LYS
1	B	157	ILE
1	B	159	LEU
1	B	320	LEU
1	B	335	GLU
1	B	345	HIS
1	B	355	SER
1	B	366	ASP
1	B	368	LYS
1	B	383	LEU
1	B	425	MET
1	B	436	LEU
1	B	466	ARG
1	B	470	ASP
1	C	18	ASP
1	C	19	ASP
1	C	74	ARG
1	C	103	ASP
1	C	110	LYS
1	C	114	GLN
1	C	137	GLU
1	C	142	VAL
1	C	152	SER
1	C	155	LYS
1	C	157	ILE
1	C	320	LEU
1	C	345	HIS
1	C	351	ARG
1	C	359	ARG
1	C	366	ASP
1	C	386	LEU
1	C	466	ARG
1	C	470	ASP
1	C	471	ASP
1	C	472	VAL

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Mol	Chain	Res	Type
1	D	19	ASP
1	D	74	ARG
1	D	110	LYS
1	D	114	GLN
1	D	117	VAL
1	D	137	GLU
1	D	152	SER
1	D	155	LYS
1	D	157	ILE
1	D	345	HIS
1	D	346	ILE
1	D	351	ARG
1	D	359	ARG
1	D	419	MET
1	D	425	MET
1	D	438	ARG
1	D	466	ARG
1	D	470	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	404	GLN
1	A	463	HIS
1	B	404	GLN
1	B	463	HIS
1	C	404	GLN
1	C	428	GLN
1	C	463	HIS
1	D	404	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	OXL	A	502	2	0,5,5	0.00	-	0,6,6	0.00	-
3	OXL	B	502	2	0,5,5	0.00	-	0,6,6	0.00	-
3	OXL	C	502	2	0,5,5	0.00	-	0,6,6	0.00	-
3	OXL	D	502	2	0,5,5	0.00	-	0,6,6	0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	OXL	A	502	2	-	0/0/4/4	0/0/0/0
3	OXL	B	502	2	-	0/0/4/4	0/0/0/0
3	OXL	C	502	2	-	0/0/4/4	0/0/0/0
3	OXL	D	502	2	-	0/0/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	464/475 (97%)	-0.08	6 (1%) 77 73	41, 65, 112, 139	0
1	B	464/475 (97%)	-0.01	10 (2%) 62 56	41, 65, 114, 140	0
1	C	464/475 (97%)	-0.02	8 (1%) 70 65	41, 65, 113, 143	0
1	D	464/475 (97%)	-0.01	7 (1%) 74 69	42, 65, 113, 137	0
All	All	1856/1900 (97%)	-0.03	31 (1%) 70 65	41, 65, 114, 143	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	472	VAL	7.2
1	A	472	VAL	5.6
1	D	472	VAL	5.6
1	B	472	VAL	5.3
1	C	75	LEU	3.3
1	B	418	LYS	3.1
1	D	470	ASP	2.6
1	B	151	VAL	2.6
1	C	139	ASP	2.5
1	C	141	VAL	2.4
1	B	141	VAL	2.4
1	D	139	ASP	2.4
1	A	136	VAL	2.3
1	B	419	MET	2.3
1	D	107	THR	2.3
1	B	135	ALA	2.2
1	B	438	ARG	2.2
1	B	420	GLN	2.2
1	D	348	ARG	2.2
1	A	85	TRP	2.2
1	B	422	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	95	VAL	2.1
1	D	471	ASP	2.1
1	A	422	THR	2.1
1	C	419	MET	2.1
1	C	418	LYS	2.1
1	C	85	TRP	2.1
1	D	418	LYS	2.0
1	C	151	VAL	2.0
1	B	93	ILE	2.0
1	A	139	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	OXL	C	502	6/6	0.97	0.19	1.54	53,60,78,79	0
3	OXL	A	502	6/6	0.96	0.18	0.82	53,62,74,78	0
3	OXL	B	502	6/6	0.97	0.16	0.27	56,61,73,80	0
3	OXL	D	502	6/6	0.96	0.15	0.16	54,64,83,86	0
2	MG	A	501	1/1	0.99	0.13	-1.07	53,53,53,53	0
2	MG	D	501	1/1	0.97	0.12	-1.47	53,53,53,53	0
2	MG	B	501	1/1	0.98	0.10	-4.33	53,53,53,53	0
2	MG	C	501	1/1	0.98	0.08	-5.66	55,55,55,55	0

6.5 Other polymers

There are no such residues in this entry.